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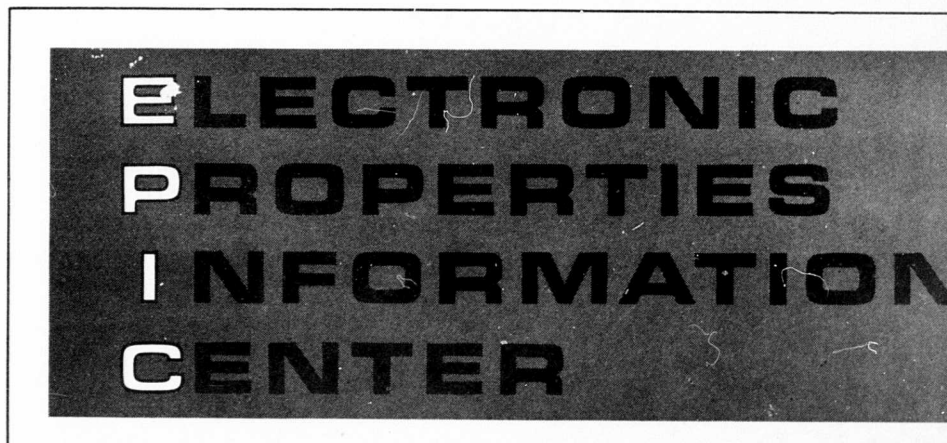
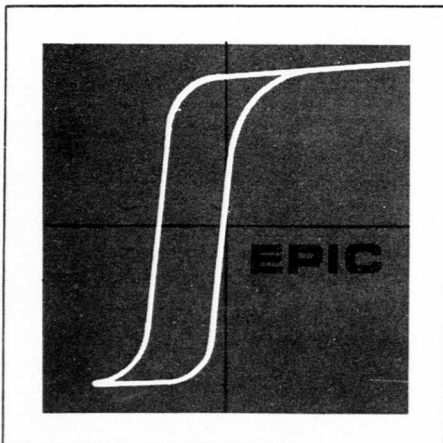
AS AD NO. 421964

# ZINC SELENIDE

Data Sheets

M. Neuberger

DS-132  
September 1963



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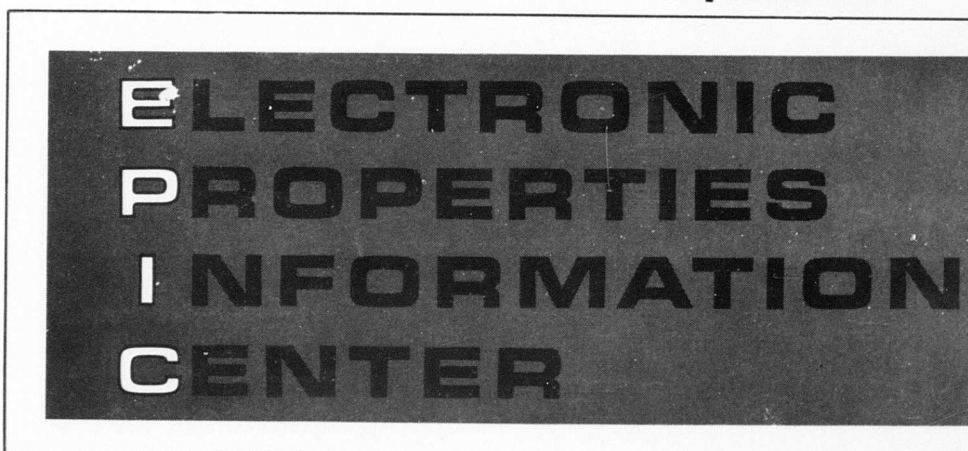
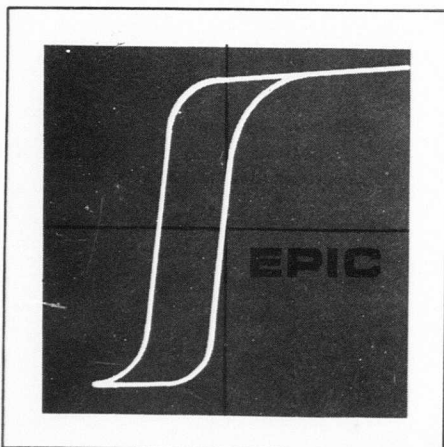
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ZINC SELENIDE

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## FOREWORD

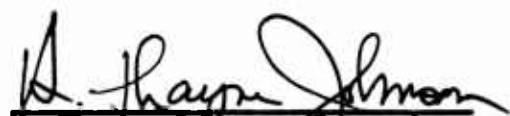
This report was prepared by Hughes Aircraft Company under Contract No. AF 33(616)-8438. The contract was initiated under Project No. 7381, Task No. 738103. The work was administered under the direction of the Directorate of Materials and Processes, Aeronautical Systems Division, with Mr. R.F. Klinger acting as Project Engineer.

Many persons have contributed to the program which this report represents. The author wishes especially to acknowledge the contributions of the following: J.J. Anders, J.W. Atwood, C.L. Blocher, D.L. Grigsby, J.J. Grossman, F.S. Harter, D.H. Johnson, H.T. Johnson, J.T. Milek, and E. Schafer.

## ABSTRACT

The Electronic Properties Information Center has been established to collect, index and abstract the literature on the electrical and electronic properties of materials and to evaluate and compile the experimental data from that literature. A modified coordinate index to the literature is machine stored and printed for manual use. The Center publishes data sheets, summary reports, thesauri, glossaries, and similar publications as sufficient information is evaluated and compiled. This report consists of the compiled data sheets on Zinc Selenide.

This report has been reviewed and is approved for publication.



H. Thayne Johnson, Supervisor  
Electronic Properties Information Center



John W. Atwood  
Project Manager

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## INTRODUCTION

In June 1961, a program was initiated under the direction of the Air Force to collect, index and abstract the literature on the electrical and electronic properties of materials and to evaluate and compile the experimental data from that literature. Placed at Hughes Aircraft Company in Culver City, California, the program, now called the Electronic Properties Information Center, was originally intended to cover ten major categories of materials: Semiconductors, Insulators, Ceramics, Ferroelectrics, Metals, Ferrites, Electroluminescent Materials, Ferromagnetics, Thermionic Emitters, and Superconductors.

During the first year, studies were completed on the Semiconductor and Insulator categories; and Ceramics was discontinued as a separate category and subsumed under the other nine. Vocabulary studies have now been completed on all categories, and retrospective documentation is virtually complete for Semiconductors and Insulators. A full index to the literature is maintained; and publications such as data sheets, summary reviews, glossaries, and thesauri are periodically issued. The use of the Center and these publications are available to anyone wishing information within the scope of the Center's objectives. A full list of publications to date appears at the end of this report.

This report contains data sheets on Zinc Selenide. The data sheets have been compiled direct from the literature. Articles are allowed to accumulate in the system until it is judged that a sufficient number are available on one material for adequate evaluation. The manual



modified coordinate index is then used to retrieve all literature on the material to be compiled. Bibliographies are checked to make sure that valuable and relevant literature is not overlooked. Then the assembled literature is given to the specialist doing the evaluation and compilation.

Evaluation is confined to primary source data except when only secondary citations are available. If equally valid data are available from several sources, all are given. Data are rejected when judged questionable because of faulty or dubious measurements, unknown sample composition, or if more reliable data are available from another source. Selection of data is based upon that which is judged most representative, precise, reliable, and covers the widest range of variables. The addition of new data to a previously evaluated property requires a reappraisal of the reported values. Older data may be deleted if the new data are judged more accurate or representative.

After a thorough analysis and evaluation, the data are compiled into data sheets which present it in its most optimum form. This will be, primarily, but not limited to, curves or tabular form. Where possible, graphs are adapted directly from the original sources. If this is not possible, they are drawn from data compiled from the articles. Where thought important, notes are entered with each graph to help the user. The references, from which the data are drawn, are shown by reference number below each graph with the full bibliographic information at the end of the data sheets. The bibliography is referred to and listed in

the order of entry into the Center (accession number). This provides a quick cross reference into the index used with the literature.

# DATA SHEET

ELECTRICAL AND ELECTRONIC PROPERTIES

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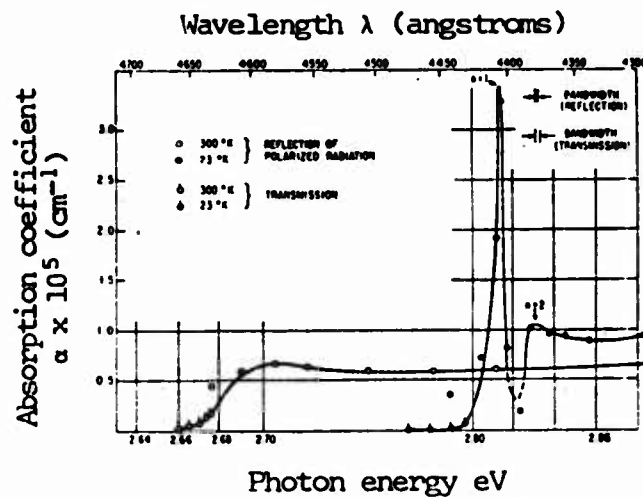
## SEMICONDUCTOR MATERIALS

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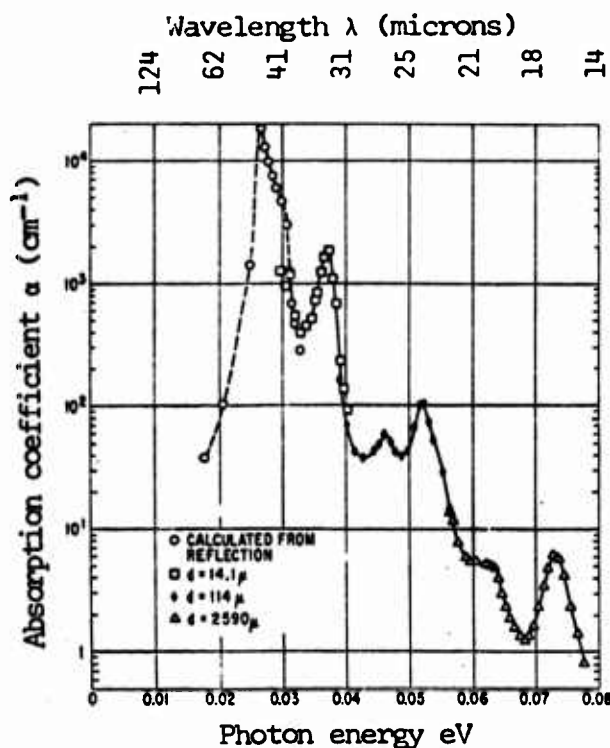
### ZINC SELENIDE

#### Absorption

Absorption coefficient for single crystal zinc selenide in the absorption edge region as a function of photon energy at 300°K and 23°K.



[Ref. 2618]



Absorption coefficient for single crystal zinc selenide in the infrared region as a function of photon energy at 300°K.

[Ref. 2618]

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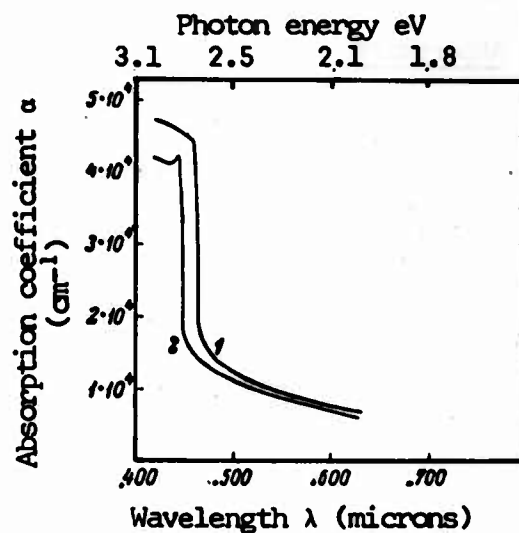
## SEMICONDUCTOR MATERIALS

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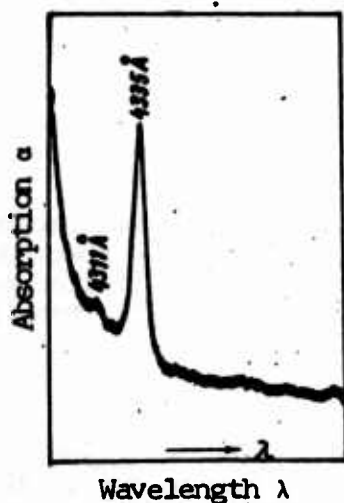
### ZINC SELENIDE

#### Absorption

Spectral distribution curves for the absorption coefficient of intrinsic, polycrystalline zinc selenide films, 0.2-2 $\mu$  thick. 1) at 393°K; 2) at 77°K.



[Ref. 690]



A microphotometer trace of the spectrum for absorption as a function of wavelength of single crystal, hexagonal zinc oxide. The sample is rotated in the extraordinary ray.

Absorption edge for a 10 $\mu$  thick sample

$\alpha = .4356\mu$  for polarized light normal to c-axis

$\alpha = .4292\mu$  for polarized light parallel to c-axis

[Ref. 5942]

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### ZINC SELENIDE

#### Absorption

Positions and polarizations of the A, B, and C exciton lines found in the spectra of single crystal, hexagonal zinc selenide at 4.2°K.

	Position	Polarization
A	4335 Å	$E_{\perp c}$
B	4311 Å	$E_{\perp c}$
C	4237 Å	$E_{\parallel c}$

Magnitudes of the valence band splitting energies  $E_{AB}$  and  $E_{AC}$ .

$E_{AB}$	$E_{AC}$
0.016eV	0.066eV

[Ref. 5942]

Wavelengths of spectral lines in edge emission spectrum of single crystal, cubic zinc selenide at 4.2°K and 77°K, type I crystals.\*

Wavelengths at 4.2°K (Å)	Wavelengths at 77°K (Å)
$L_1$ —4410	$N_1$ —4421
$L_2$ —4420	$N_2$ —4436
$L_2'$ —4424	
$L_3$ —4448	
$L_3'$ —4454	
$L_{3a}$ —4490	
$L_{3b}$ —4551	
$L_{3c}$ —4602	
$L_{3d}$ —4654	
$L_{3e}$ —4712	

Wavelengths at 4.2°K (Å)	Wavelengths at 77°K (Å)
$L_1$ —4420	$N_1$ —4388
$L_2$ —4427	$N_2$ —4590
$L_3$ —4437	$N_{3a}$ —4638
$L_4$ —4453	
$L_5$ —4488	
$L_6$ —4506	
$L_7$ —4512	
$L_8$ —4548	
$L_9$ —4558	
$L_{10}$ —4598	
$L_{10a}$ —4655	
$L_{10b}$ —4712	
$L_{10c}$ —4770	
$L_{10d}$ —4832	

Wavelengths of spectral lines in edge emission spectrum of single crystal, cubic zinc selenide at 4.2°K and 77°K, type II crystals.\*

\*The two types of crystals may be due to different host lattice defects.

[Ref. 2500]

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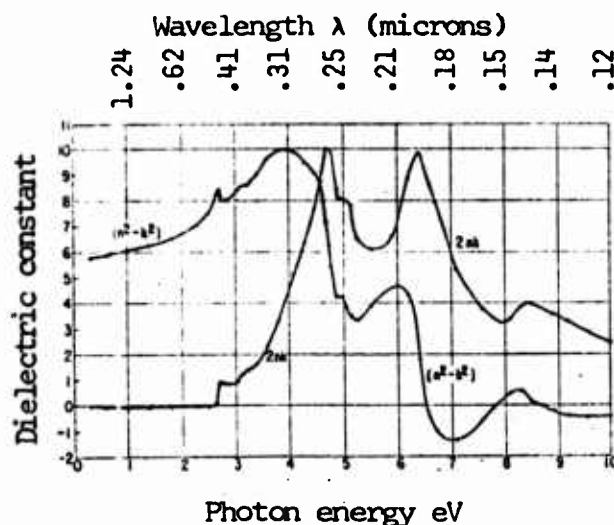
### ZINC SELENIDE

#### Debye Temperature

Symbol	Value	Type	Temperature	Ref.
$\theta_D$	400°K	polycrystalline	80°K	3030

#### Dielectric Constant

Symbol	Value	Type	Temperature	Ref.
$\epsilon_0$	9.1	single crystal, cubic	298°K	10288
$\epsilon_0$	$8.1 \pm 0.3$	single crystal	300°K	2618
$\epsilon_\infty$	$5.75 \pm 0.1$	single crystal	300°K	2618



Real and imaginary parts of the dielectric constant,  $(n^2-k^2)$  and  $2nk$ , respectively, for single crystal zinc selenide as a function of photon energy at 300°K.  $n$  = refractive index;  $k$  = extinction coefficient;  $n$  and  $k$  are calculated from optical data.

[Ref. 2618]

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### ZINC SELENIDE

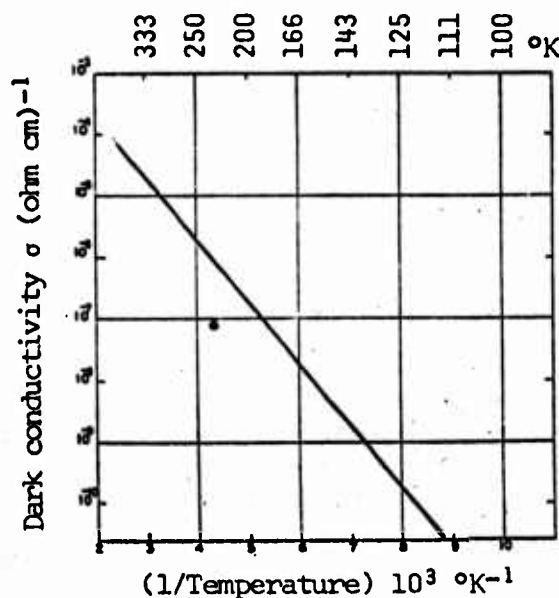
#### Effective Mass

Symbol	Value	Type	Test Conditions
$m_n^*$	$\sim 0.1 m_0$	single crystal, n- and p-type	optical measurement, 300°K
$m_p^*$	$\sim 0.6 m_0$	single crystal, n- and p-type	optical measurement, 300°K [Ref. 2618]

#### Electrical Conductivity

Symbol	Value (ohm cm) <sup>-1</sup>	Dopant	Temperature	Test Conditions
$\sigma$	$4 \times 10^{-2}$	bromine, arsenic	300°K	{ single crystal, dark conductivity
$\sigma$	$7 \times 10^{-6}$	bromine, antimony	300°K	

[Ref. 780]



Dark conductivity as a function of temperature for single crystal zinc selenide doped with bromine and antimony.

[Ref. 780]

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### ZINC SELENIDE

#### Electrical Resistivity

Symbol	Value	Type	Temperature	Ref.
$\rho$	$10^8$ to $10^9$ ohm cm	0.2 to $2\mu$ thick films polycrystalline, pure	300°K	690

#### Electroacoustic Properties

Symbol	Value	Type	Temperature		Ref.
LO	0.031 eV	single crystal	300°K	optical data	2618
TO	0.026 eV				
LO	0.03 eV		4.2°K	optical data	2500
TO	0.027 eV				

#### Energy Bands

Symbol	Value	Temperature	Test Conditions
dE <sub>g</sub> /dP	$6.0 \times 10^{-6}$ eV/atm	300°K	single crystal, pure, experimental wave- length, $\lambda = .481\mu$ [Ref. 273]
	$0.49 \times 10^{-6}$ eV/atm (maximum shift)	300°K	
	$-2.0 \times 10^{-6}$ eV/atm	300°K	
dE <sub>g</sub> /dT	$7.2 \times 10^{-4}$ eV/°K	90 - 400°K	single crystal, pure [Ref. 826]



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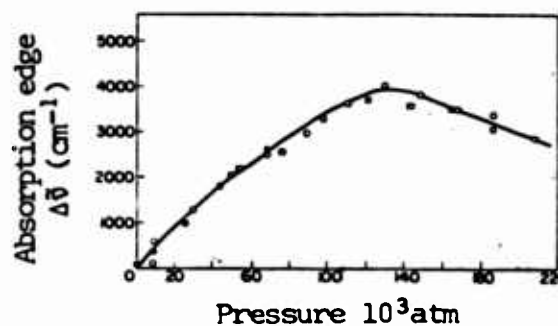
## SEMICONDUCTOR MATERIALS

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### ZINC SELENIDE

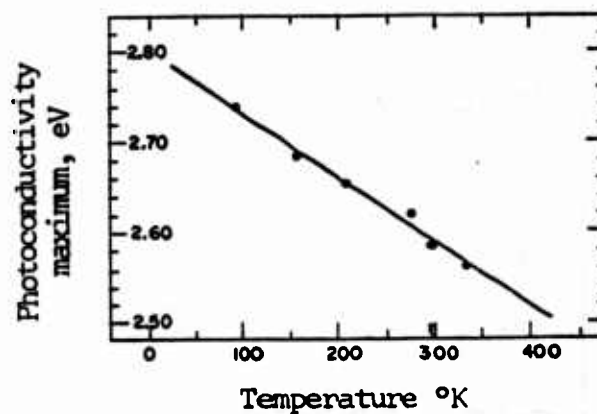
#### Energy Bands

Shift of zinc selenide absorption edge with pressure ( $\nu_0 = 20,800 \text{ cm}^{-1}$ ,  $\alpha = 65 \text{ cm}^{-1}$ ).  
Samples are single crystal, pure.



[Ref. 273]

The location of the photoconductivity maximum (or maxima) as a function of temperature for single crystal zinc selenide.



[Ref. 826]

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## SEMICONDUCTOR MATERIALS

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### ZINC SELENIDE

#### Energy Gap

Symbol	Value	Type	Temperature	Ref.
$E_g$	2.9 (calc.)	polycrystalline film,	0°K	10250
	2.59	zincblende symmetry,	295°K	
	2.47	1 $\mu$ film, optical-	398°K	
	2.38	absorption measurement	496°K	
$E_g$	2.67	single crystal	297°K	5201
$E_g$	2.6	polycrystalline film	300°K	7174
$E_g$	2.786	polycrystalline film	77.3°K	634
$E_g$	2.81 $\pm$ 0.01	single crystal electro-optical	4°K	2618
$E_g$	2.83	cubic	4.2°K	2500

#### Energy Levels

Symbol	Value (eV)	Dopant	Type	Ref.
$E_D$	0.21	Bromine	single crystal	780
$E_A$	0.6	Copper	single crystal	780
$E_A$	0.6	Silver	single crystal	780
$E_A$	0.7	Antimony	single crystal	780

Energy gaps calculated from reflectance data. Double entries correspond to peaks split by spin-orbit interaction. Sample is single crystal. Temperature = 300°K.

Energy transitions	$L_1' \rightarrow L_1$	$L_1' \rightarrow L_2$	$X_1' \rightarrow X_1$ $X_1' \rightarrow X_1$	$X_1' \rightarrow X_1(?)$
	4.9	9.1		
	5.3	9.6	6.7	8.5

[Ref. 5949]

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### ZINC SELENIDE

#### Energy Levels

Transitions at the L-absorption edge showing spin orbit splitting of the  $L_3$  valence band in thin film zinc selenide, cubic sample.

Value (eV)	Temperature (°K)
4.94	20
4.92	78
4.89	200
4.80	297

[Ref. 2850]

Exciton absorption peaks associated with valence and conduction bands at  $\Gamma$  for single crystal zinc selenide at 4°K. Principal peak at  $2.81 \pm 0.01$  eV.

	1st. peak at $\Gamma$	2nd. peak at L
Exciton transition	2.7 eV	4.75 eV
Interband transition	3.15 eV	5.1 eV
Spin orbit valence-band split	0.45 eV	0.35 eV

[Ref. 2618]

#### Mobility

Symbol	Value	Type	Ref.
$\mu_n$	530 $\text{cm}^2(\text{Vsec})^{-1}$	single crystal, n-type, $\rho=1$ ohm cm	300°K 5954
$\mu_n$	$\sim 100$ $\text{cm}^2(\text{Vsec})^{-1}$		300°K 2911

Material	Temp. range (°C)	Mobility ( $\text{cm}^2/\text{v-sec}$ )		Carrier concentration ( $\text{cm}^{-3}$ )	
		Electrons	Holes	27°K	200°K
ZnSe:Cu	200-400		11	...	$10^{10}$
ZnSe:Cu, Se-fired	130-260		16	...	$10^{11}$
ZnSe, Se-fired	200-400		15	...	$10^{10}$
ZnSe:Ga	27-400	80		$10^8$	...
ZnSe:Ga, Zn-fired	27	150		$10^{18}$	...
ZnSe, Zn-fired	27-250	260		$10^{18}$	...

Mobility and carrier concentration for single crystal zinc selenide, variously doped at temperatures from 300°K to 673°K.

[Ref. 2618]

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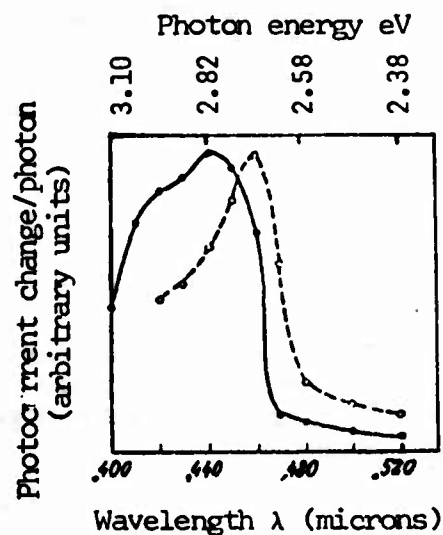
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### ZINC SELENIDE

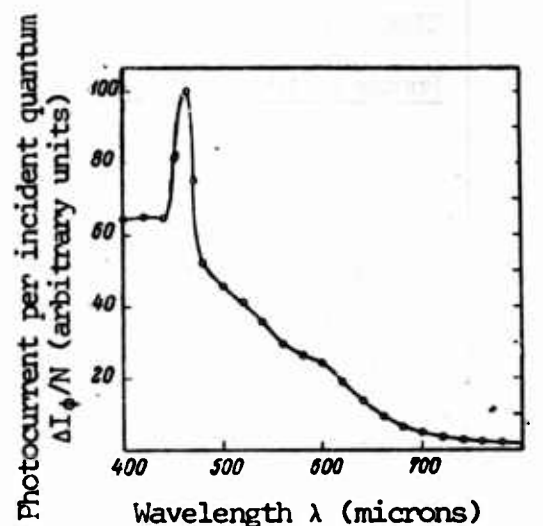
#### Photoelectronic Properties

Spectral distribution curves for the photoconductivity of intrinsic, polycrystalline zinc selenide films, 0.2 - 2 $\mu$  thick at 300°K.

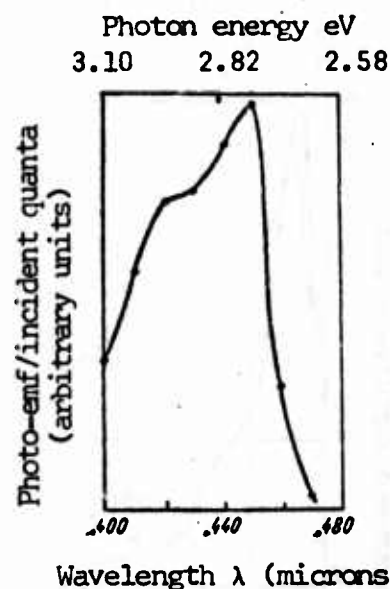


Decrease in electrical conductivity on illumination in a polycrystalline zinc selenide film. The broken curve indicates photoconductivity.

[Ref. 622]



[Ref. 690]



The photo-emf response spectrum of a polycrystalline zinc selenide sample.

[Ref. 622]

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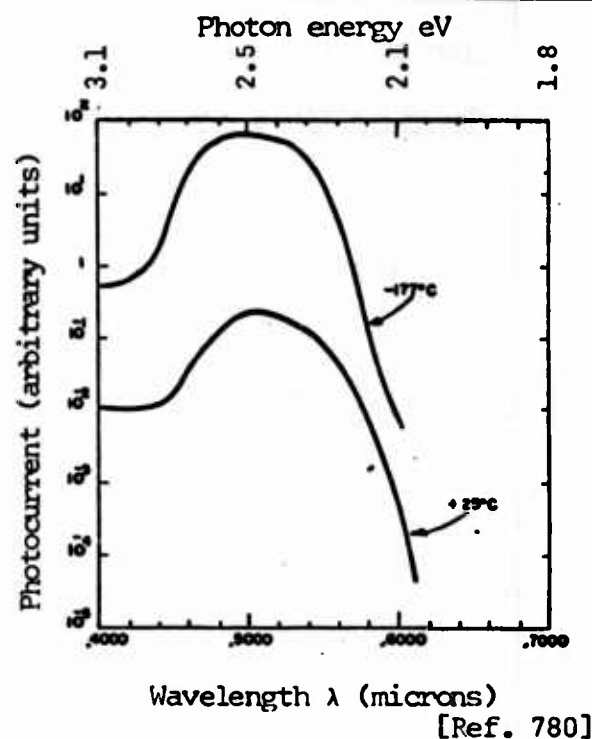
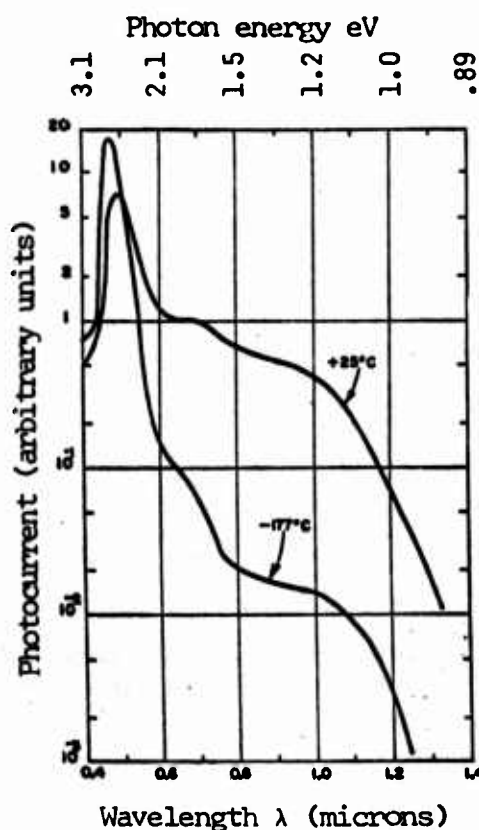
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### ZINC SELENIDE

#### Photoelectronic Properties

Spectral response curves for photoconductivity in single crystal, bromine-silver-doped zinc selenide.



Spectral response curves for photoconductivity in single crystal, bromine-antimony-doped zinc selenide.

[Ref. 780]

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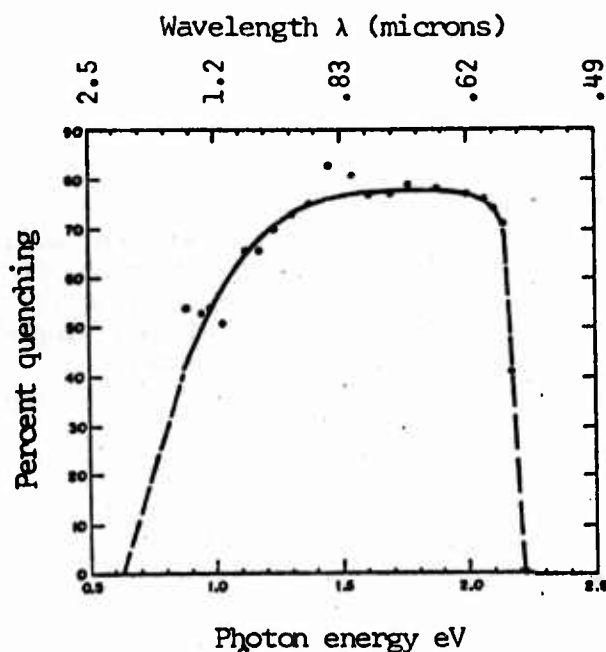
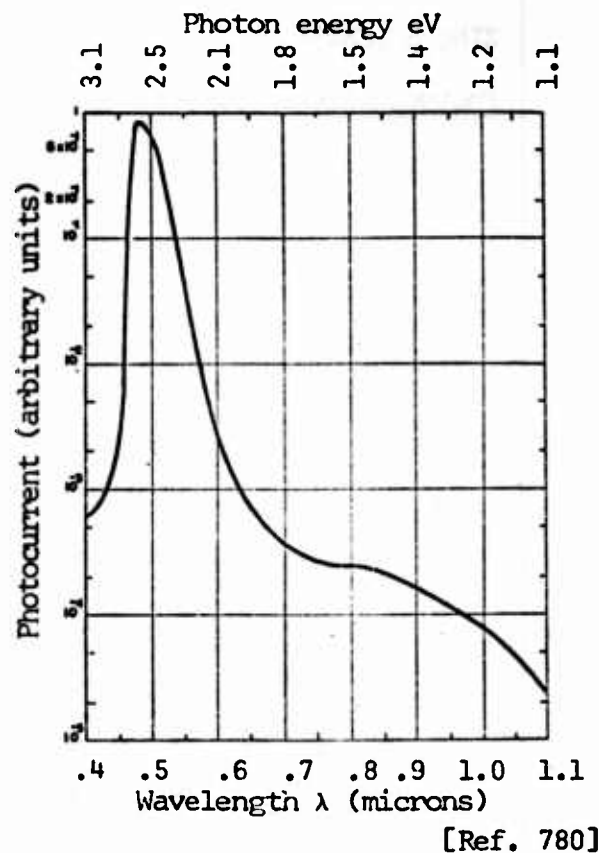
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### ZINC SELENIDE

#### Photoelectronic Properties

Spectral response curve for photoconductivity in single crystal, bromine-arsenic-doped zinc selenide.



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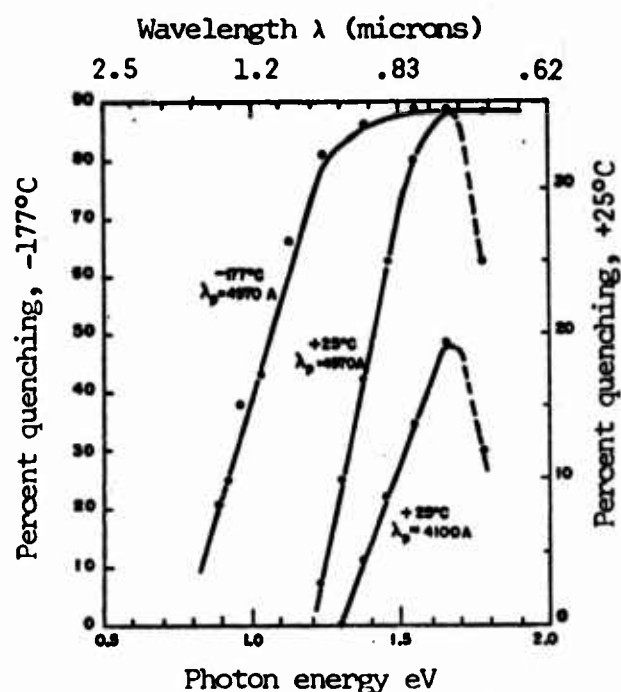
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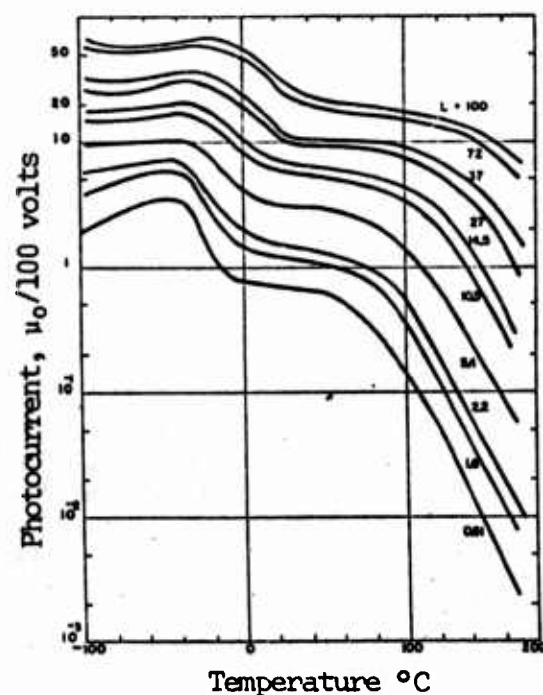
### ZINC SELENIDE

#### Photoelectronic Properties

Infrared quenching spectra for single crystal, bromine-antimony-doped zinc selenide.



[Ref. 780]



Variation of photocurrent with temperature for different photo-intensities of excitation for bromine-antimony-doped, single crystal zinc selenide.

[Ref. 780]

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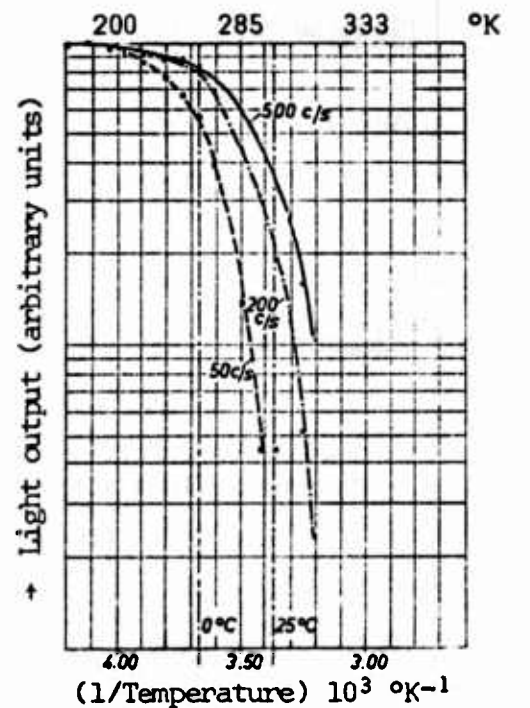
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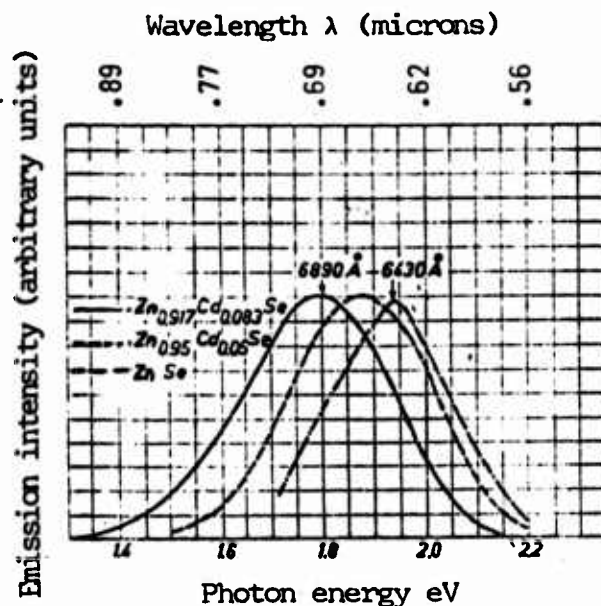
### ZINC SELENIDE

#### Photon Electroluminescence

Photon Electroluminescence: light output as a function of temperature for a copper-aluminum-doped zinc selenide red phosphor.



[Ref. 6458]



Emission intensity as a function of wavelength for a copper-aluminum-doped zinc selenide red phosphor and a mixed zinc cadmium selenide phosphor.

[Ref. 6458]



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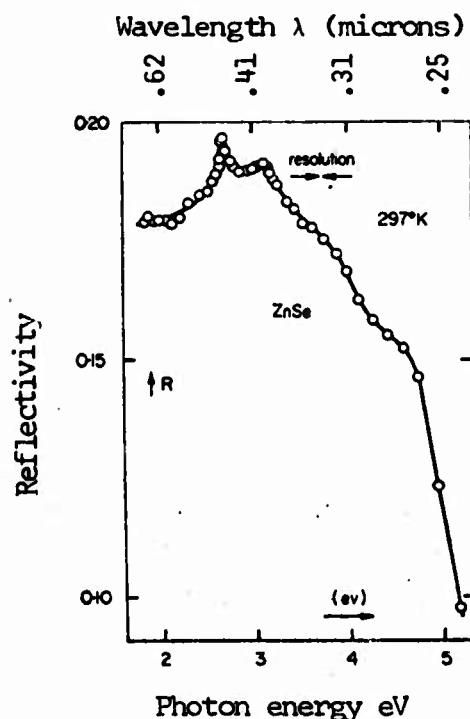
### ZINC SELENIDE

#### Piezoelectric Properties

Symbol	Value	Type	Temperature	Ref.
$d_{31}$	$0.31 \times 10^{-12}$	single crystal, cubic	298°K	10288
$d_{14}$	$1.10 \times 10^{-12}$			

#### Reflectivity

	Type	Temperature
Maximum reflectivity at $.4242\mu$	single crystal, hexagonal	4.2°K
Minimum reflectivity at $.4232\mu$		
(polarized light parallel to c-axis)		
		[Ref. 5942]



Reflectivity spectrum of single crystal, cubic zinc selenide as a function of wavelength at 297°K.

[Ref. 3935]

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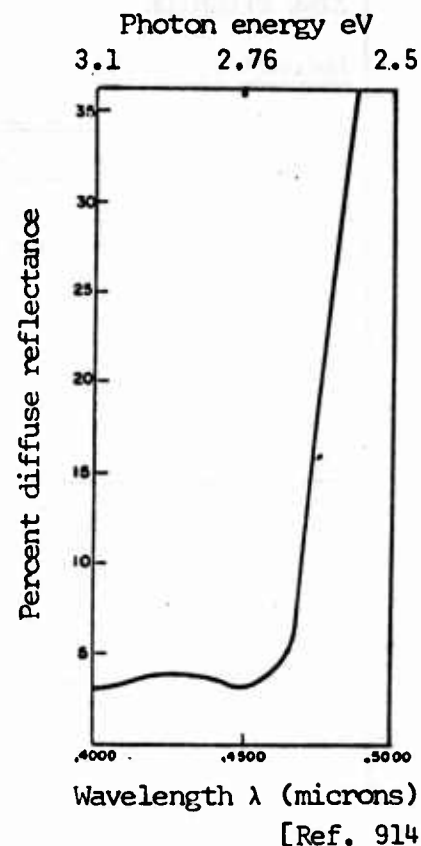
## SEMICONDUCTOR MATERIALS

September 1963

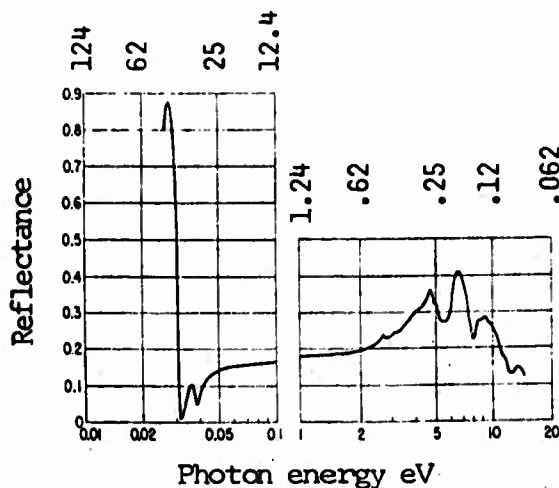
### ZINC SELENIDE

#### Reflectivity

Diffuse reflectance spectra of polycrystalline zinc selenide at 300°K.



Wavelength λ (microns)



Reflectance of single crystal zinc selenide for near-normal incidence as a function of photon energy at 300°K.

[Ref. 2618]

# DATA SHEET

ELECTRICAL AND ELECTRONIC PROPERTIES

MATERIALS CENTRAL  
AERONAUTICAL SYSTEMS DIVISION  
AIR FORCE SYSTEMS COMMAND

## SEMICONDUCTOR MATERIALS

September 1963

### ZINC SELENIDE

#### Refractive Index

Symbol	Value	Ref.
n	2.89 $\lambda = .589\mu$	7359

#### Thermal Conductivity

Symbol	Value	Temperature	Ref.
k	$33 \times 10^3$ cal/cm sec deg	300°K	636
k	$30-33 \times 10^3$ cal/cm sec deg		3477

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